

10/705, 966

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: ssspta1204bxsd

PASSWORD :

TERMINAL (ENTER 1, 2, 3, OR ?):2

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 12:00:35 ON 23 DEC 2004

```
=> fil regf  
'REGF' IS NOT A VALID FILE NAME  
SESSION CONTINUES IN FILE 'HOME'  
Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files  
that are available. If you have requested multiple files, you can
```

specify a corrected file name or you can enter "IGNORE" to continue accessing the remaining file names entered.

=> fil reg
COST IN U.S. DOLLARS

SINCE FILE ENTRY	TOTAL SESSION
0.21	0.21

FILE 'REGISTRY' ENTERED AT 17:01:03 ON 23 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 22 DEC 2004 **HIGHEST RN** 802006-11-7
DICTIONARY FILE UPDATES: 22 DEC 2004 **HIGHEST RN** 802006-11-7

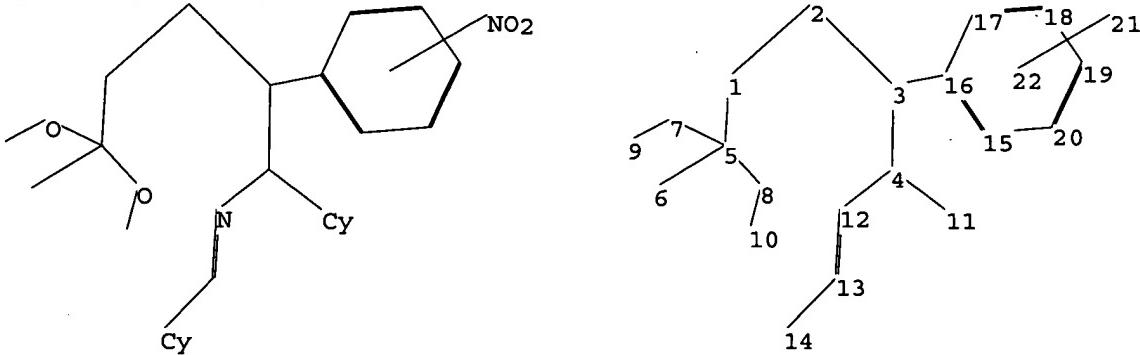
TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>
Uploading C:\Program Files\Stnexp\Queries\10705466.str



```

chain nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 21
ring nodes :
15 16 17 18 19 20
chain bonds :
1-2 1-5 2-3 3-4 3-16 4-11 4-12 5-6 5-7 5-8 7-9 8-10 12-13 13-14
ring bonds :
15-16 15-20 16-17 17-18 18-19 19-20
exact/norm bonds :
4-11 4-12 5-7 5-8 7-9 8-10 12-13 13-14
exact bonds :
1-2 1-5 2-3 3-4 3-16 5-6
normalized bonds :
15-16 15-20 16-17 17-18 18-19 19-20

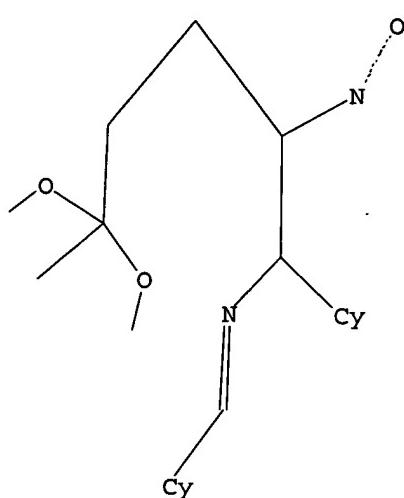
```

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS
10:CLASS 11:Atom 12:CLASS 13:CLASS 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom
19:Atom 20:Atom 21:CLASS 22:CLASS

L1 STRUCTURE UPLOADED

=> d query

L1



Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 17:01:17 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s 11 full
FULL SEARCH INITIATED 17:01:22 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED 2 ITERATIONS 2 ANSWERS
SEARCH TIME: 00.00.01

L3 2 SEA SSS FUL L1

=> fil caplus
COST IN U.S. DOLLARS

SINCE FILE ENTRY	TOTAL SESSION
------------------	---------------

FULL ESTIMATED COST

155.42

155.63

FILE 'CAPLUS' ENTERED AT 17:01:25 ON 23 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 23 Dec 2004 VOL 141 ISS 26
FILE LAST UPDATED: 22 Dec 2004 (20041222/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

```
=> s 13
L4          2 L3
=> d 14 abs ibib
```

L4 ANSWER 1 OF 2 CAPIUS COPYRIGHT 2004 ACS on STN
AB This document discloses a process for preparing a pure cis isomer from a mixture of cis-trans isomers of formula $RC(OMe)2CH_2CH_2CH(NO_2)CH(Ar)N:CHAR$ (I) [Ar = (un)substituted Ph, etc.; R = alkyl] comprising the steps of :
(a) dispersing a mixture of cis- and trans-I in an inert solvent wherein said cis isomer is less soluble than said trans isomer; (b) heating said dispersion to completely dissolve said trans isomer; (c) maintaining said heating step to allow interconversion of said cis and trans isomer; (d) cooling said mixture thereby crystallizing the cis isomer; (e)

separating said crystalline cis isomer from said solvent. Cis isomers of formula I are useful intermediates in the synthesis of cis isomers of benzamide piperidine compds. which exhibit activity as NK-1 receptor antagonists.

ACCESSION NUMBER: 2004428898 CAPIUS

DOCUMENT NUMBER: 141:6912

TITLE: Process for converting a cis-trans mixture of substituted benzylidene amines into the pure cis isomer

INVENTOR(S): Humphrey, John Michael; Tom, Norma Jacqueline

PATENT ASSIGNEE(S): Pfizer Products Inc., USA

SOURCE: PCT Int. Appl. 21 pp.

CODEN: PIKKDZ

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004043908	A1	20040527	WO 2003-IB4953	20031103
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MN, MM, MX, MZ, NI, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: BW, GH, GM, KE, LS, MM, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TU, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD,				

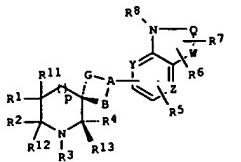
TG US 2004254380 A1 20041216 US 2003-705466 20031110
PRIORITY APPN. INFO.: US 2002-425946P P 20021112

OTHER SOURCE(S): CASREACT 141:6912; MARPAT 141:6912

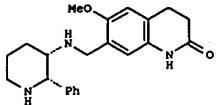
REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE

FORMAT

=> d 14 2 abs ibib



I



II

AB Title compds. I [Q = C:NH, C:CH₂, C:S, C:O, SO, SO₂; A = CH, CH₂, C(alkyl), CH(alkyl), C(CF₃), or CH(CF₃) with the proviso that when B is present, A = CH, C(alkyl), or C(CF₃); B = absent, CH₂, or ethylene; Y, Z = N, CH, provided that both are not N; G = NH(CH₂)q, S(CH₂)q, O(CH₂)q; q = 0-1 with the proviso that when q = 0, G = NH₂, SH, OH; W = 1-3 carbon linking group, including spiro assemblies; p = 0-2; R₃ = H, acyl, carboxy, Ph, heterocyclic, alkyl, etc.; R₁, R₂, R₁₁-R₁₃ = H, alkyl, etc., or R₁₂-R₁₃ together with the carbon atoms to which they are attached form a 5- or 6-membered heterocyclic ring, etc.; R₄ = Ph, pyridyl, thiényl, etc.; R₅-R₈ = H, alkyl, S(O)₁₋₂-alkyl, S(O)₁₋₂-aryl, alkoxy, halo, Ph, etc.] were prepared. Approx. 100 synthetic examples and over 100 precursor preps. were provided. For instance, 4-aminophenol was acylated with 3-chloropropionyl chloride (CH₂Cl₂, H₂O, NaHCO₃, room temperature, 4 h) and the product treated with AlCl₃ at 210°C for 10 min effecting cyclization to the hydroxy quinolone intermediate. The intermediate was O-methylated (acetone, Me₂SO₄, K₂CO₃, room temperature, 16 h) and formylated in the 7 position (CH₂Cl₂, AlCl₃, Cl₂CHOMe) to give 7-formyl-6-methoxy-1H-1,2,3,4-tetrahydroquinolin-2-one. Reductive alkylation of the quinolone with (2S,3S)-3-amino-2-phenylpiperidine (a. PhMe, 3Å mol. sieves; b. dichloroethane, NaHB(OAc)₃, room temperature, 16 h) yielded II. Compds. I are NK-1 receptor antagonists, i.e., substance P receptor antagonists. At least one stereoisomer of the example compds. had a binding affinity, as measured by

L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2004 ACS on STN (Continued)
Ki, of at least 600 nM. I are used in the treatment and prevention of a wide variety of central nervous system disorders, inflammatory disorders, cardiovascular disorders, ophthalmic disorders, etc.

ACCESSION NUMBER: 2001:762988 CAPLUS
DOCUMENT NUMBER: 135:331346
TITLE: Synthesis of benzamide piperidine containing compounds as substance P antagonists
INVENTOR(S): Arnold, Eric Platt; Chappie, Thomas Allen; Huang, Jianhua; Humphrey, John Michael; Nagel, Arthur Adam; O'Neill, Brian Thomas; Sobolov-Jayne, Susan Beth; Vincent, Lawrence Albert
PATENT ASSIGNEE(S): Pfizer Products Inc., USA
SOURCE: PCT Int'l Appl., 209 pp.
CODEN: PIXXD2

DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001077100	A2	20011018	WO 2001-IB629	20010406
WO 2001077100	A3	20020307		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BE, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KE, LC, LK, LR, LS, LT, LV, MA, MD, MG, MN, MM, MK, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AA, AZ, BY, KG, KZ, MD, RO, TO, IN, RW: GH, GM, KE, LS, MM, ME, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GM, ML, MR, NE, SN, TD, TG				
US 2003087925	A1	20030508	US 2001-811218	20010316
CA 2405089	AA	20011018	CA 2001-2405089	20010406
EP 1272484	A2	20030108	EP 2001-919702	20010406
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IE, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
BR 2001009936	A	20030506	BR 2001-9936	20010406
JP 2004501072	T2	20040115	JP 2001-575573	20010406
EE 20020588	A	20040415	EE 2002-588	20010406
NZ 521346	A	20040730	NZ 2001-521346	20010406
BG 107135	A	20030630	BG 2002-107135	20020923
ZA 2002008072	A	20031008	ZA 2002-8072	20021008
NO 2002004874	A	20021118	NO 2002-4874	20021009
PRIORITY APPLN. INFO.: US 2000-195922P				P 20000410
OTHER SOURCE(S): MARPAT 135:331346				

=> fil reg		SINCE FILE	TOTAL
COST IN U.S. DOLLARS		ENTRY	SESSION
FULL ESTIMATED COST		5.98	161.61
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)		SINCE FILE	TOTAL
CA SUBSCRIBER PRICE		ENTRY	SESSION
		-1.40	-1.40

FILE 'REGISTRY' ENTERED AT 17:02:31 ON 23 DEC 2004
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
 provided by InfoChem.

STRUCTURE FILE UPDATES: 22 DEC 2004 HIGHEST RN 802006-11-7
 DICTIONARY FILE UPDATES: 22 DEC 2004 HIGHEST RN 802006-11-7

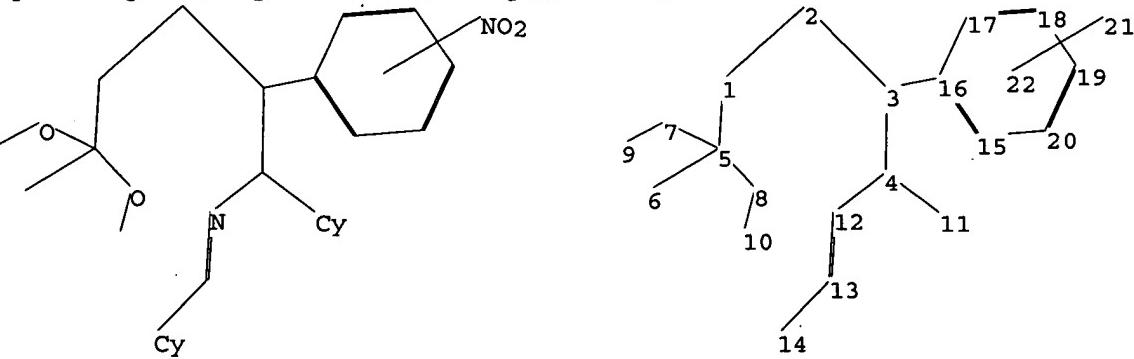
TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
 information enter HELP PROP at an arrow prompt in the file or refer
 to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>
 Uploading C:\Program Files\Stnexp\Queries\10705466.str



```

graph TD
    2 --- 1
    1 --- 5
    5 --- 9
    5 --- 7
    5 --- 8
    5 --- 10
    9 --- 6
    9 --- 1
    7 --- 6
    7 --- 8
    3 --- 16
    16 --- 17
    16 --- 18
    16 --- 19
    16 --- 22
    17 --- 18
    18 --- 21
    19 --- 20
    20 --- 15
    15 --- 12
    12 --- 13
    12 --- 11
    13 --- 14
    4 --- 12
    4 --- 11
    11 --- 10
    10 --- 8
    8 --- 1
    3 --- 4
    4 --- 16
    16 --- 17
    17 --- 18
    18 --- 19
    19 --- 20
    15 --- 16
    15 --- 20
    16 --- 17
    17 --- 18
    18 --- 19
    19 --- 20
    12 --- 13
    13 --- 14
    14 --- 1
  
```

chain nodes :
 1 2 3 4 5 6 7 8 9 10 11 12 13 14 21
 ring nodes :
 15 16 17 18 19 20
 chain bonds :
 1-2 1-5 2-3 3-4 3-16 4-11 4-12 5-6 5-7 5-8 7-9 8-10 12-13 13-14
 ring bonds :
 15-16 15-20 16-17 17-18 18-19 19-20
 exact/norm bonds :
 4-11 4-12 5-7 5-8 7-9 8-10 12-13 13-14
 exact bonds :

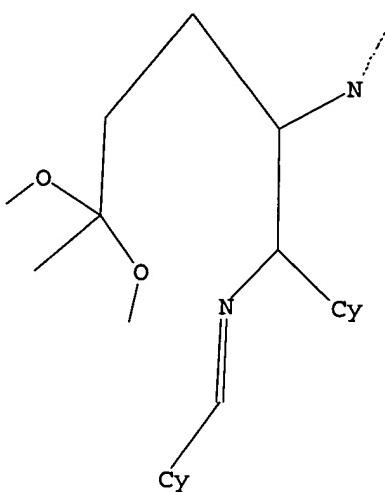
1-2 1-5 2-3 3-4 3-16 5-6
normalized bonds :
15-16 15-20 16-17 17-18 18-19 19-20

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS
10:CLASS 11:Atom 12:CLASS 13:CLASS 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom
19:Atom 20:Atom 21:CLASS 22:CLASS

L5 STRUCTURE UPLOADED

=> d query

L5 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 15
SAMPLE SEARCH INITIATED 17:03:26 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 1 TO ITERATE

100.0% PROCESSED 1 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 1 TO 80
PROJECTED ANSWERS: 0 TO 0

L6 0 SEA SSS SAM L5

=> s 15 full
FULL SEARCH INITIATED 17:03:30 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 22 TO ITERATE

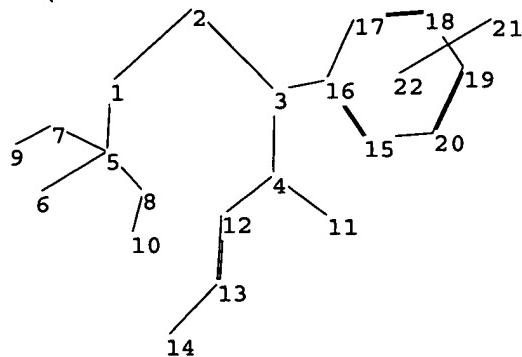
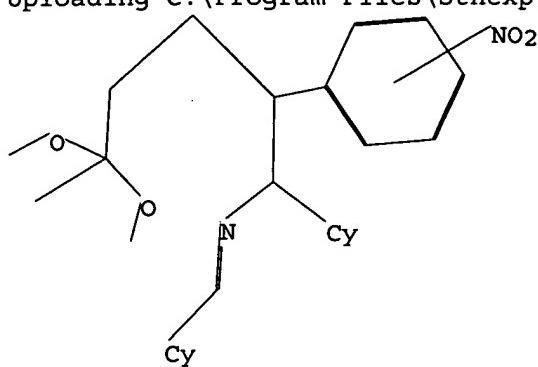
100.0% PROCESSED 22 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L7 0 SEA SSS FUL L5

=>

Uploading C:\Program Files\Stnexp\Queries\10705466.str



chain nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 21

ring nodes :

15 16 17 18 19 20

chain bonds :

1-2 1-5 2-3 3-4 3-16 4-11 4-12 5-6 5-7 5-8 7-9 8-10 12-13 13-14

ring bonds :

15-16 15-20 16-17 17-18 18-19 19-20

exact/norm bonds :

4-11 4-12 5-7 5-8 7-9 8-10 12-13 13-14

exact bonds :

1-2 1-5 2-3 3-4 3-16 5-6

normalized bonds :

15-16 15-20 16-17 17-18 18-19 19-20

Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS

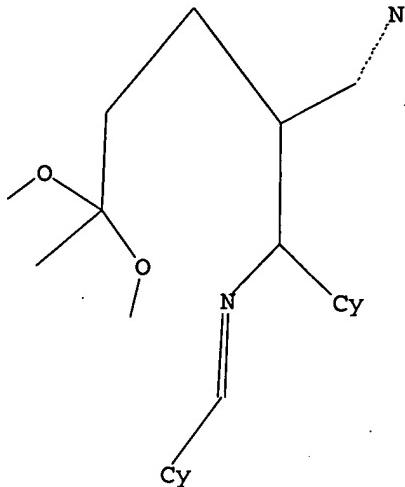
10:CLASS 11:Atom 12:CLASS 13:CLASS 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom

19:Atom 20:Atom 21:CLASS 22:CLASS

L8 STRUCTURE UPLOADED

=> d query

L8 STR



Structure attributes must be viewed using STN Express query preparation.

```
=> s 18
SAMPLE SEARCH INITIATED 17:04:10 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -      11 TO ITERATE

100.0% PROCESSED      11 ITERATIONS          0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE   **COMPLETE**
                        BATCH    **COMPLETE**
PROJECTED ITERATIONS:    22 TO      418
PROJECTED ANSWERS:        0 TO       0

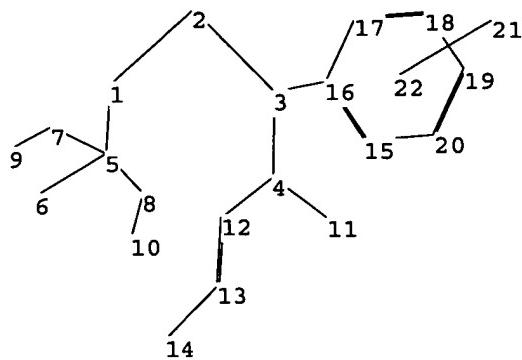
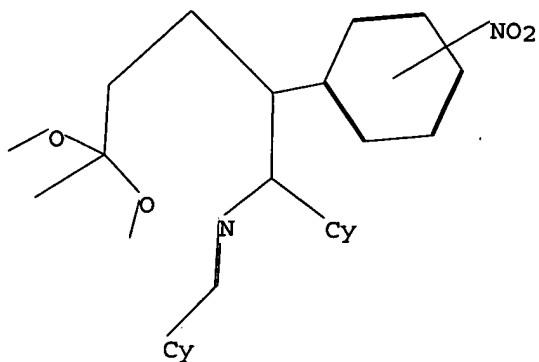
L9      0 SEA SSS SAM L8

=> s 18 full
FULL SEARCH INITIATED 17:04:14 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -    111 TO ITERATE

100.0% PROCESSED      111 ITERATIONS         0 ANSWERS
SEARCH TIME: 00.00.01

L10     0 SEA SSS FUL L8

=>
Uploading C:\Program Files\Stnexp\Queries\10705466.str
```



chain nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 21

ring nodes :

15 16 17 18 19 20

chain bonds :

1-2 1-5 2-3 3-4 3-16 4-11 4-12 5-6 5-7 5-8 7-9 8-10 12-13 13-14

ring bonds :

15-16 15-20 16-17 17-18 18-19 19-20

exact/norm bonds :

4-11 4-12 5-7 5-8 7-9 8-10 12-13 13-14

exact bonds :

1-2 1-5 2-3 3-4 3-16 5-6

normalized bonds :

15-16 15-20 16-17 17-18 18-19 19-20

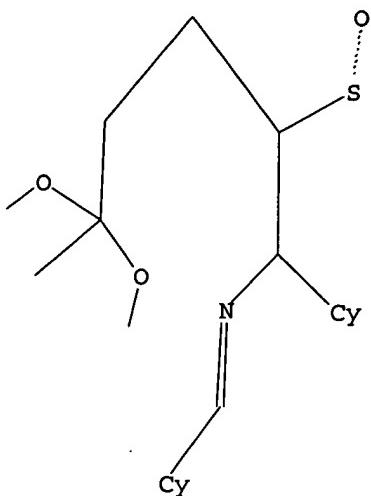
Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS
 10:CLASS 11:Atom 12:CLASS 13:CLASS 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom
 19:Atom 20:Atom 21:CLASS 22:CLASS

L11 STRUCTURE UPLOADED

=> d query

L11 STR



Structure attributes must be viewed using STN Express query preparation.

```
=> s l11
SAMPLE SEARCH INITIATED 17:05:53 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -      1 TO ITERATE

100.0% PROCESSED      1 ITERATIONS          0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:    1 TO      80
PROJECTED ANSWERS:       0 TO      0

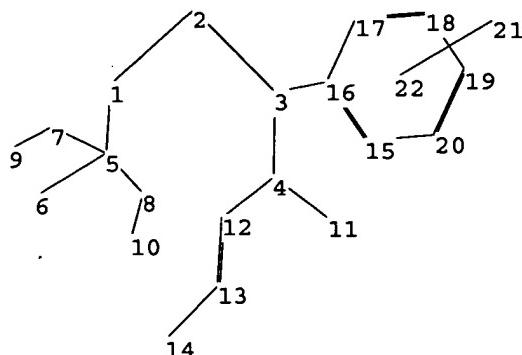
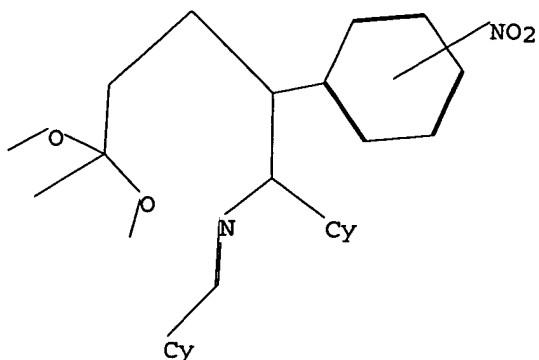
L12      0 SEA SSS SAM L11

=> s l11 full
FULL SEARCH INITIATED 17:05:57 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -      4 TO ITERATE

100.0% PROCESSED      4 ITERATIONS          0 ANSWERS
SEARCH TIME: 00.00.01

L13      0 SEA SSS FUL L11

=>
Uploading C:\Program Files\Stnexp\Queries\10705466.str
```



chain nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 21

ring nodes :

15 16 17 18 19 20

chain bonds :

1-2 1-5 2-3 3-4 3-16 4-11 4-12 5-6 5-7 5-8 7-9 8-10 12-13 13-14

ring bonds :

15-16 15-20 16-17 17-18 18-19 19-20

exact/norm bonds :

4-11 4-12 5-7 5-8 7-9 8-10 12-13 13-14

exact bonds :

1-2 1-5 2-3 3-4 3-16 5-6

normalized bonds :

15-16 15-20 16-17 17-18 18-19 19-20

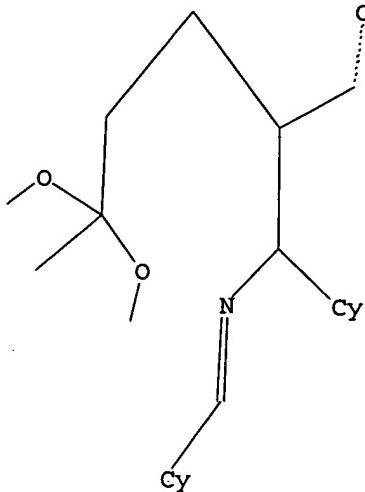
Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS
 10:CLASS 11:Atom 12:CLASS 13:CLASS 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom
 19:Atom 20:Atom 21:CLASS 22:CLASS

L14 STRUCTURE UPLOADED

=> d query

L14 STR



Structure attributes must be viewed using STN Express query preparation.

```
=> s l14
SAMPLE SEARCH INITIATED 17:06:43 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -      10 TO ITERATE
```

100.0% PROCESSED	10 ITERATIONS	0 ANSWERS
SEARCH TIME: 00.00.01		

FULL FILE PROJECTIONS:	ONLINE	**COMPLETE**
	BATCH	**COMPLETE**
PROJECTED ITERATIONS:	11 TO	389
PROJECTED ANSWERS:	0 TO	0

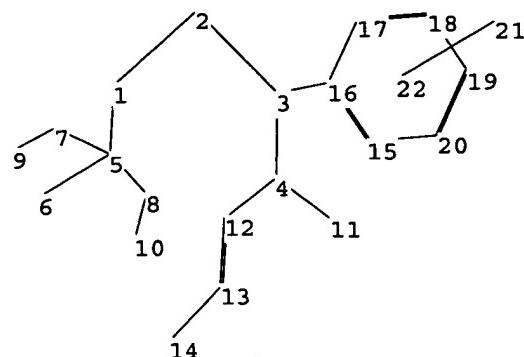
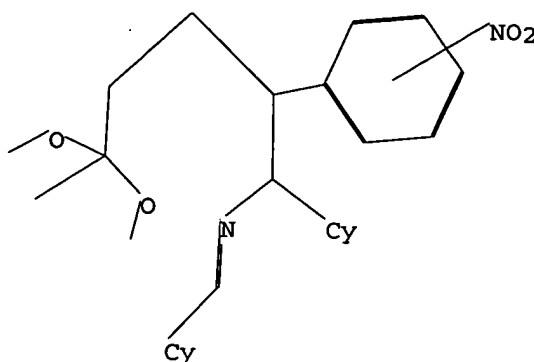
L15 0 SEA SSS SAM L14

```
=> s l14 full
FULL SEARCH INITIATED 17:06:51 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -      275 TO ITERATE
```

100.0% PROCESSED	275 ITERATIONS	0 ANSWERS
SEARCH TIME: 00.00.01		

L16 0 SEA SSS FUL L14

```
=>
Uploading C:\Program Files\Stnexp\Queries\10705466.str
```



chain nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 21

ring nodes :

15 16 17 18 19 20

chain bonds :

1-2 1-5 2-3 3-4 3-16 4-11 4-12 5-6 5-7 5-8 7-9 8-10 12-13 13-14

ring bonds :

15-16 15-20 16-17 17-18 18-19 19-20

exact/norm bonds :

4-11 4-12 5-7 5-8 7-9 8-10 12-13 13-14

exact bonds :

1-2 1-5 2-3 3-4 3-16 5-6

normalized bonds :

15-16 15-20 16-17 17-18 18-19 19-20

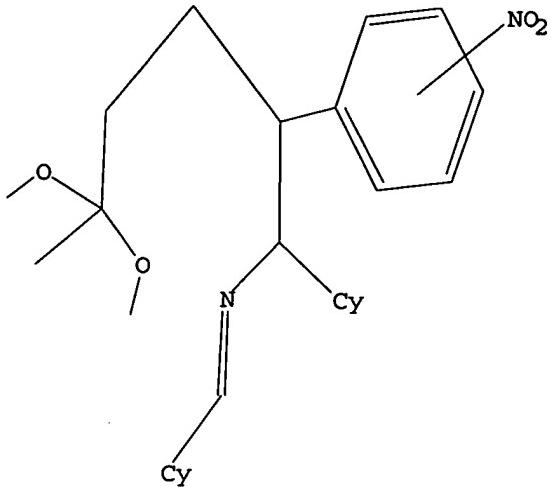
Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS
 10:CLASS 11:Atom 12:CLASS 13:CLASS 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom
 19:Atom 20:Atom 21:CLASS 22:CLASS

L17 STRUCTURE UPLOADED

=> d query

L17 STR



Structure attributes must be viewed using STN Express query preparation.

```
=> s 117
SAMPLE SEARCH INITIATED 17:08:06 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -      0 TO ITERATE
```

100.0% PROCESSED	0 ITERATIONS	0 ANSWERS
SEARCH TIME: 00.00.01		

FULL FILE PROJECTIONS:	ONLINE	**COMPLETE**
	BATCH	**COMPLETE**
PROJECTED ITERATIONS:	0 TO	0
PROJECTED ANSWERS:	0 TO	0

L18 0 SEA SSS SAM L17

```
=> s 117 full
FULL SEARCH INITIATED 17:08:10 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -      1 TO ITERATE
```

100.0% PROCESSED	1 ITERATIONS	0 ANSWERS
SEARCH TIME: 00.00.01		

L19 0 SEA SSS FUL L17

=> logoff y		
COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	784.24	945.85
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.40

STN INTERNATIONAL LOGOFF AT 17:15:35 ON 23 DEC 2004